Trying 3106016892...Open

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LOGINID:ssspta1204rxw
PASSWORD:
TERMINAL (ENTER 1, 2, 3, OR ?):2

NEWS 1 Web Page URLs for STN Seminar Schedule - N. America
NEWS 2 Dec 17 The CA Lexicon available in the CAPLUS and CA files
NEWS 3 Feb 06 Engineering Information Encompass files have new names
NEWS 4 Feb 16 TOXLINE no longer being updated
NEWS 5 Apr 23 Search Derwent WPINDEX by chemical structure
NEWS 6 Apr 23 PRE-1967 REFERENCES NOW SEARCHABLE IN CAPLUS AND CA
NEWS 7 May 07 DGENE Reload
NEWS 8 Jun 20 Published patent applications (A1) are now in USPATFULL
NEWS 9 JUL 13 New SDI alert frequency now available in Derwent's
DWPI and DPCI

NEWS EXPRESS

July 11 CURRENT WINDOWS VERSION IS V6.0b,

CURRENT MACINTOSH VERSION IS V5.0C (ENG) AND V5.0JB (JP),

AND CURRENT DISCOVER FILE IS DATED 06 APRIL 2001

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STN Operating Hours Plus Help Desk Availability

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FILE 'HOME' ENTERED AT 09:34:50 ON 30 JUL 2001

=> file reg

COST IN U.S. DOLLARS
SINCE FILE TOTAL
ENTRY SESSION
FULL ESTIMATED COST
0.15
0.15

FILE 'REGISTRY' ENTERED AT 09:34:55 ON 30 JUL 2001 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

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STRUCTURE FILE UPDATES: 29 JUL 2001 HIGHEST RN 349446-89-5 DICTIONARY FILE UPDATES: 29 JUL 2001 HIGHEST RN 349446-89-5

TSCA INFORMATION NOW CURRENT THROUGH January 11, 2001

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Structure search limits have been increased. See HELP SLIMIT for details.

=>Testing the current file.... screen

ENTER SCREEN EXPRESSION OR (END):end

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Uploading c:\stnexp4\queries\09652376.str

L1 STRUCTURE UPLOADED

=> que L1

L2 QUE L1

=> d

L2 HAS NO ANSWERS L1 STR

Structure attributes must be viewed using STN Express query preparation. L2 $\,$ OUE $\,$ L1 $\,$

=> s 12

SAMPLE SEARCH INITIATED 09:35:32 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 446 TO ITERATE

100.0% PROCESSED 446 ITERATIONS SEARCH TIME: 00.00.01

23 ANSWERS

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS: PROJECTED ANSWERS:

7654 TO 10186 173 TO 747

L3 23

23 SEA SSS SAM L1

=> d scan

L3 23 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN Benzenepropanoic acid, .alpha.-(1,3-benzodioxol-5-ylhydrazono)-2,4,5trimethoxy-, methyl ester (9CI)

MF C20 H22 N2 O7

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HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):0

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FULL SEARCH INITIATED 09:35:57 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 9102 TO ITERATE

100.0% PROCESSED 9102 ITERATIONS

1 ANSWERS

SEARCH TIME: 00.00.01

1.4

1 SEA CSS FUL L1

=> d

L4 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2001 ACS

RN 6906-65-6 REGISTRY

CN Benzene, 1,2,4-trimethoxy-5-propyl- (7CI, 8CI, 9CI) (CA INDEX NAME) OTHER NAMES:

CN 1,2,4-Trimethoxy-5-propylbenzene

CN Dihydroasarone

FS 3D CONCORD

MF C12 H18 O3

LC STN Files: BEILSTEIN*, CA, CAOLD, CAPLUS, CASREACT, TOXLINE, TOXLIT (*File contains numerically searchable property data)

09/652,376

- 5 REFERENCES IN FILE CA (1967 TO DATE)
- 5 REFERENCES IN FILE CAPLUS (1967 TO DATE)
- 1 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

=> file caplus caold

COST IN U.S. DOLLARS

SINCE FILE TOTAL SESSION ENTRY 135.37 135.52

FULL ESTIMATED COST

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=> s 14

L5 6 L4

=> d 1-6 bib ab hitstr

- L_5 ANSWER 1 OF 6 CAPLUS COPYRIGHT 2001 ACS
- 1995:740149 CAPLUS AN
- DN 123:187711
- TI Studies on crude drugs effective on visural larva migrans. XVI. Mobility inhibition and nematocidal activity of asarone and related phenylpropanoids on second-stage larvae of Toxocara canis
- Sugimoto, Naoki; Goto, Yoshihisa; Akao, Nobuaki; Kiuchi, Fumiyuki; Kondo, ΑU Kaoru; Tsuda, Yoshisuke
- Sch. Med., Kanazawa Univ., Kanazawa, 920, Japan
- Biol. Pharm. Bull. (1995), 18(4), 605-9 CODEN: BPBLEO; ISSN: 0918-6158
 - Journal
- DΤ LΑ English
- AB The in vitro effect of asarone, the nematocidal principle of the rhizome of Acorus calamus, on second-stage larvae of Toxocara canis is composed of

two independent actions: one is a fast acting inhibition of the larval mobility and the other is a slow acting larvicidal action. Mobility of the larvae was rapidly inhibited when they were incubated with asarone. Dye exclusion assay revealed that larvae were alive at this stage, and their mobility was restored after the first inhibition, suggesting that this effect was temporary and reversible. However, when the mobility

decreased again during prolonged incubation, the cellular viability of larvae disappeared, showing that they were killed by the compd. above

two-stage effect of asarone was almost identical in two geometrical isomers ((E)- and (Z)-asarone). Di- and tri-methoxypropenyl or propylbenzenes carrying two methoxy groups at a vicinal position on a benzene ring showed, more a less, a two-stage effect of this type. These two actions were suggested to be separable by an appropriate modification of the structure.

TΤ 6906-65-6

RL: BAC (Biological activity or effector, except adverse); PRP (Properties); BIOL (Biological study)

(studies on crude drugs effective on visural larva migrans. XVI. Mobility inhibition and nematocidal activity of asarone and related phenylpropanoids on second-stage larvae of Toxocara canis)

6906-65-6 CAPLUS RN

Benzene, 1,2,4-trimethoxy-5-propyl- (7CI, 8CI, 9CI) (CA INDEX NAME) CN

L5 ANSWER 2 OF 6 CAPLUS COPYRIGHT 2001 ACS

ΑN 1990:138847 CAPLUS

DN 112:138847

ΤI Potential antipsychotic agents. 5. Synthesis and antidopaminergic properties of substituted 5,6-dimethoxysalicylamides and related compounds

ΑU Hoegberg, Thomas; Bengtsson, Stefan; De Paulis, Tomas; Johansson, Lars; Stroem, Peter; Hall, Haakan; Oegren, Sven Ove

CS CNS Res. Dev., Astra Res. Cent. AB, Soedertaelje, S-151 85, Swed.

J. Med. Chem. (1990), 33(4), 1155-63 CODEN: JMCMAR; ISSN: 0022-2623

DT /Journal

LA English

OS CASREACT 112:138847

AΒ A series of 3-substituted 5,6-dimethoxysalicylamides (I) have been synthesized from the corresponding 2,5,6-trimethoxybenzoic acids. Relaxation times T1 and carbon chem. shifts of the methoxy groups in I showed that the 6-methoxy group adopts a nearly perpendicular orientation and the 5-methoxy group takes on a more coplanar orientation with respect to the ring plane in soln. The salicylamides I display a very high and stereoselective affinity for the [3H]spiperone and [3H]raclopride binding sites in vitro. Regioisomers of salicylamides I also exhibit pronounced, but lower than I, affinity for the [3H]spiperone binding site. The structural requirements were further assessed by studies of the related amino analogs and hydroxy analog. 3-Bromo compd. II (FLB 463) was

studied

in various in vivo models and compared with the dopamine-D2 antagonists sulpiride, raclopride, eticlopride, and haloperidol. The high potency of I to selectively block dopamine-D2 receptors in vitro and in vivo

with indications on a low potential for motor side effects makes it a very

interesting new member of the class of substituted salicylamides.

IT 6906-65-6P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation) (prepn. and carboxylation of)

RN 6906-65-6 CAPLUS

CN Benzene, 1,2,4-trimethoxy-5-propyl- (7CI, 8CI, 9CI) (CA INDEX NAME)

L5 ANSWER 3 OF 6 CAPLUS COPYRIGHT 2001 ACS

AN 1983:517832 CAPLUS

DN 99:117832

TI Repellency and toxicity of 55 insect repellents to red-winged blackbirds (Agelaius phoeniceus)

AU Schafer, E. W., Jr.; Jacobson, M.

CS Denver Wildl. Res. Cent., USFWS, Denver, CO, 80225, USA

SO J. Environ. Sci. Health, Part A (1983), A18(4), 493-502

CODEN: JESEDU; ISSN: 0360-1226

DT Journal

LA English

AB A joint research program was initiated in 1979 to investigate the potential avian repellency and toxicity of 55 selected insect repellents originating from or related to naturally occurring chems. Seven of the chems. or exts. tested exhibited avian repellency and 2 of these were considered to be moderately active, with 50% avian repellency concns. of 0.237 (trans-asarone [2883-98-9]) and 0.240% (safrole [94-59-7]). None of the 55 chem. or exts. exhibited acute oral toxicity at .ltoreq.100 mg/kg to the red-winged blackbird (Agelaius phoeniceus).

IT 6906-65-6

RL: ADV (Adverse effect, including toxicity); BIOL (Biological study) (repellency and toxicity of, to red-winged blackbirds)

RN 6906-65-6 CAPLUS

CN Benzene, 1,2,4-trimethoxy-5-propyl- (7CI, 8CI, 9CI) (CA INDEX NAME)

L5 ANSWER 4 OF 6 CAPLUS COPYRIGHT 2001 ACS

AN 1982:100901 CAPLUS

DN 96:100901

TI Constituents of Acorus calamus: structure of acoramone. Carbon-13 NMR spectra of cis- and trans-asarone

AU Patra, Amarendra; Mitra, Alok K.

CS Dep. Chem., Univ. Coll. Sci., Calcutta, 700009, India

J. Nat. Prod. (1981), 44(6), 668-9

CODEN: JNPRDF; ISSN: 0163-3864

101.7615

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oxa compds. - (VI)

6906-65-6 CAOLD

6906-65-6

El'tsov, A. V.; Minkin, V. I.; Tsereteli, I. Yu.

DT Journal LΑ English ΑB The phenylpropane derivs. isoeugenol Me ether, .gamma.-asarone, cis-asarone, trans-asarone, and the new natural product acoramone (I) were isolated along with asarylaldehyde from the oil of A. calamus. 13C-NMR signal assignments of cis- and trans-asarone and their dihydroderivative are also reported. IT 6906-65-6 RL: BOC (Biological occurrence); BIOL (Biological study); OCCU (Occurrence) (of Acorus calamus) 6906-65-6 CAPLUS RN CN Benzene, 1,2,4-trimethoxy-5-propyl- (7CI, 8CI, 9CI) (CA INDEX NAME) Pr-n MeO. MeO OMe ANSWER 5 OF 6 CAPLUS COPYRIGHT 2001 ACS L5 ΑN 1969:31614 CAPLUS DN 70:31614 ΤI Isolation of 2,4,5-trimethoxyallylbenzene from Caesulia axillaries oil ΑU Devgan, O. N.; Bokadia, M. M. csر Southern Illinois Univ., Carbondale, Ill., USA SO Aust. J. Chem. (1968), 21(12), 3001-3 CODEN: AJCHAS Journal DТ English LΑ AΒ The phenolic ether obtained from the essential oil of C. axillaries has been shown to be 2,4,5-trimethoxyallyl-benzene, on the basis of chem. and spectroscopic evidence. It has tentatively been named as .gamma.-asarone. IT · 6906-65-6P RL: PREP (Preparation) (from Caesulia axillaris oil) RN 6906-65-6 CAPLUS CN Benzene, 1,2,4-trimethoxy-5-propyl- (7CI, 8CI, 9CI) (CA INDEX NAME) Pr-n Me0 MeO OMe 1/L5 ANSWER 6 OF 6 CAOLD COPYRIGHT 2001 ACS AN CA65:8851d CAOLD

CN Benzene, 1,2,4-trimethoxy-5-propyl- (7CI, 8CI, 9CI) (CA INDEX NAME)

=> log y

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	24.04	159.56
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                to PHARMASEARCH
NEWS EXPRESS August 15 CURRENT WINDOWS VERSION IS V6.0c,
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